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Attention: Mr. Warren M. Cheek

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February 13, 2009

VIA FACSIMILE
(Total Page: 6)

Your Ref: 2006-0369A
Our Ref: 546533
U.S. Patent Application No. 10/572,639
Applicants: Yoshitaka TOMIGAHARA et al.

ACKNOWLEDGED WITH RECEIPTION
WENDEROTH, LIND & PONACK, L.L.P.
KNOX PRIORITY: ILW
DATE RECEIVED: Feb 13, 09

Dear Mr. Cheek:

Thank you for your letter of February 11, 2009 concerning the Examiners' request.

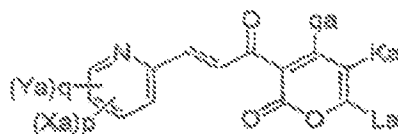
We would like to offer the following comments.

1) Ring A

The applicants' preference is, first, benzene, and then pyridine.

2) Location of N in pyridine ring

Please refer the following formula. The location shown in the formula is preferred.



3) Preference of a substituent for each Xa, Ya, La, Ka and qa

Please refer to the attached sheet. The order of preference is shown by the symbols ①, ②, ③,

4) Order of importance of Xa, Ya, La, Ka and qa

The order is Xa > qa > La > Ka > Ya. That is, first is Xa, next is qa,

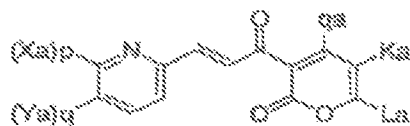
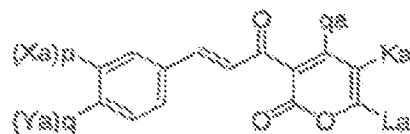
To: WENDEROTH, LIND & PONACK, L.L.P.

February 13, 2009

Page 2 of 2

5) Point of attachment of X and Y on the ring

Please refer to the following formulas.



Please respond to the Examiner.

Kindly acknowledge receipt of this letter.

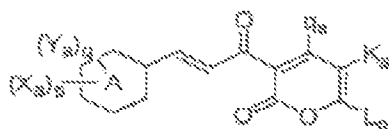
Sincerely yours,
AOYAMA & PARTNERS

Mitsuo TANAKA

MT/xme
Encl.

~~1~~

4. (Original) A cinnamoyl compound represented by the formula (IV):



(IV)

wherein

A represents a ①benzene ring or a ②pyridine ring,

X_a is a substituent on a carbon atom, and represents ⑥a C1-C10 alkyl group substituted with a cyano group; ⑦a C1-C10 alkyl group substituted with a tetrahydropyran-4-ylidene group; ⑧a C2-C10 alkenyl group substituted with a halogen atom or a cyano group; ⑨a C2-C10 alkenyl group substituted with a C1-C10 alkoxy carbonyl group; ⑩a C3-C10 alkynyl group substituted with a hydroxyl group; ⑪an a₁-x₁-b-x₁'- group (wherein a₁ represents a methyl group substituted with a C1-C10 alkylthio group, a methyl group substituted with a C1-C10 alkylsulfinyl group, a methyl group substituted with a C1-C10 alkylsulfonyl group, a C2-C10 alkenyl group, a C2-C10 alkynyl group, a x₁O-CO- group (wherein x₁ represents a C1-C10 alkyl group, or a C2-C10 alkyl group substituted with a hydroxyl group), a carboxyl group, a x₁x₁'N-CO- group (wherein x and x' are the same or different, and represent a hydrogen atom or a C1-C10 alkyl group), an a₁-NH-CO- group (wherein a₁ represents a C2-C10

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alkyl group substituted with a C1-C10 alkoxy group), an a_1' -CO- group (wherein a_1' represents a morpholino group), a rr' N-CH₂- group (wherein r and r' are as defined above), a r_0 -(O)₁-CONH-CH₂- group (wherein r_0 represents a C1-C10 alkyl group, and 1 represents 0 or 1), a r -OCH₂- group (wherein r is as defined above), a r_0 -CO- group (wherein r_0 is as defined above), a cyano group, or a sulfomethyl group, r_1 represents a C1-C10 alkylene group, r_1' represents a single bond or a C1-C10 alkylene group, and b represents an oxy group, a thio group, a sulfinyl group, a sulfonyl group, or an imino group); ⑤an a_1 -y-CO-NH- group (wherein a_1 represents a C2-C10 alkyl group substituted with a C1-C10 alkoxy group, and y represents an oxy group or an imino group); ⑥a r_0 -COCO-NH- group (wherein r_0 is as defined above); ⑦an a_1 -x-NH- group (wherein a_1 represents a C2-C10 alkenyl group, or a C1-C10 alkyl group substituted with a C1-10 alkoxy group, a C1-C10 alkoxycarbonyl group, a carboxy group or a cyano group, and x represents a carbonyl group or a sulfonyl group); ⑧an a_1 -NHCO- group (wherein a_1 represents a C1-C10 alkoxy group, or a C3-C10 alkenyloxy group, or a r_0 -SO₂- group (wherein r_0 is as defined above), or a C2-C10 alkyl group substituted with a hydroxyl group or a C1-C10 alkoxy group, or a C1-C10 alkyl group substituted with a r O-CO- group (wherein r is as defined above), a cyano group or an aminocarbonyl group, or a r O-

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CO-(xO-COCH₃)CH₂- group (wherein x is as defined above)); ⑤ an a₂-NHSO₂- group (wherein a₂ represents a C2-C10 alkyl group substituted with a C1-C10 alkoxy group); ⑥a r₂ON=CH- group (wherein r₂ is as defined above); ⑦a r₂NHCSNH- group (wherein r₂ is as defined above); ⑧a r₂NHC(-Sr₂')-N- group (wherein r₂ is as defined above, r₂' is the same as the different from r₂ and has the same meaning as r₂ has); or ⑨ a (r₂O)_pP(-O)CH₂- group (wherein r₂ is as defined above);

p represents 1, 2 or 3, and when p is 2 or more, X_ss are the same or different;

Y_s represents ①a halogen atom, ②a nitro group, ③a r₂CO-NH- group (wherein r₂ is as defined above), ④a C1-C10 alkyl group or ⑤a C1-C10 alkoxy group;

q represents 0, 1 or 2, and when q is 2 or more, Y_ss are the same or different;

q_s represents ①a x_s-O- group (wherein x_s represents a hydrogen atom, a C1-C10 alkyl group, a C3-C10 alkenyl group, a C3-C10 alkynyl group, a C1-C10 alkyl group substituted with a x_sx_s'N-CH₂- group (wherein x_s and x_s' are as defined above), a xOCH₂- group (wherein x is as defined above), a x_s-CO- group (wherein x_s is as defined above), a C1-C10 alkoxycarbonyl group, a carboxy group, an aminocarbonyl group or a cyano group, or a r₁-r₁-group (wherein r₁ represents a phenyl group or a pyridyl group, and r₁ is as defined above)); ②a piperidino group; ③a morpholino

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group; or ④a x_1x_2N - group (wherein x_1 and x_2 are the same or different, and represent a hydrogen atom, a C1-C10 alkyl group, a C3-C10 alkenyl group, a C3-C10 alkynyl group, or a C2-C10 alkyl group substituted with a C1-C10 alkoxy group, provided that x_1 and x_2 are not a hydrogen atom at the same time);

K_n represents ①a hydrogen atom, ③a halogen atom or ②a C1-C10 alkyl group, and L_n represents ②a hydrogen atom or ①a C1-C10 alkyl group; or

K_n and L_n together may form a ②C1-C10 alkylene group or a ①1,3-butadienylene group;

the term "as defined above" used for the same symbols among plural substituents means that the plural substituents independently represent the same meaning as that described above and, among the plural substituents, although the selection range of substituents to be selected is the same, selected substituents may be the same or different as long as they are selected within the range.

* * * COMMUNICATION RESULT REPORT (FEB. 13. 2009 9:40AM) * * *

FAX HEADER 1:
FAX HEADER 2:TRANSMITTED/STORED : FEB. 13. 2009 9:40AM
FILE MODE OPTION

ADDRESS

RESULT

PAGE

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E-18 88% FACSIMILE CONNECTION

1-2030333615=FAX5510

/0669490362=AOYAMA

/09-02-13-15:53/001-001



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February 13, 2009

VIA FACSIMILE
(Total Page: 6)

Your Ref: 2006-0869A
Our Ref: 846833
U.S. Patent Application No. 10/872,639
Applicants: Yoshitaka TOMIYAMA et al.

ACKNOWLEDGED WITH APPLICATION
WENDEROTH, LIND & PONACK, L.L.P.
RECEIVED BY: [Signature]
DATE RECEIVED: Feb 13, 2009

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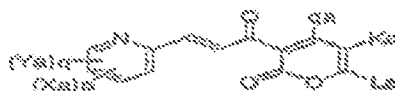
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